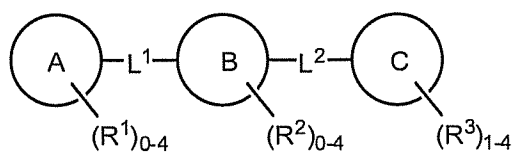


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

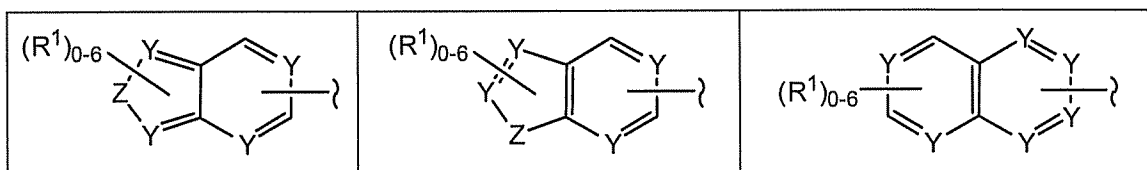
1. (currently amended) A compound for modulating c-Kit activity according to Formula I,



I

or a pharmaceutically acceptable salt, thereof, wherein,

ring A is:



wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S-, and -N(R⁷)-;

each R¹ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R¹, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R¹⁰;

L¹ is a single bond;

ring B is phenyl ~~a five- to ten-membered aryl or a five- to ten-membered heterocyclyl~~;

each R² is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R², together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R¹⁵;

L² is selected from -N(H)N(H)C(=O)N(H)-, -CH₂N(H)C(=O)N(H)-, -CH₂OC(=O)N(H)-, and -XCH₂C(=O)N(H)-; wherein X is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-; and any C-H

of L^2 is optionally $C-R^{20}$;

ring C is phenyl or pyridyl;

each R^3 is independently selected from halogen, trihalomethyl, -CN,

-NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl; provided R^3 is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide,

wherein there exists at least one of R^3 that is halogen or trihalomethyl;

~~two adjacent of R^3 ; together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{25} ;~~

R^4 is selected from -H, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two of R^4 , when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^5 is selected from -H, -CN, -NO₂, -OR⁴, -S(O)₀₋₂R⁴, -CO₂R⁴, optionally substituted C₁₋₆alkyl, optionally substituted C₁₋₆alkenyl, and optionally substituted C₁₋₆alkynyl;

R^7 is selected from -H, optionally substituted C₁₋₆alkyl, -SO₂N(R⁴)R⁴, -CO₂R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl; and

each of R^{10} , each of R^{15} , each of R^{20} , and each of R^{25} is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -CO₂R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

provided:

1) when both ring B and ring C are phenyl:

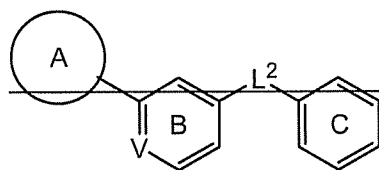
- a) and the compound comprises ring B-CH₂N(H)C(=O)N(H)-ring C, then L¹ must be a single bond; R³ can not comprise a group of the formula -O(CH₂)₂₋₄-N-piperazine that is *ortho*- to L²; and ring A cannot be a 5-methyl-[1,2,4]-oxadiazol-3-yl radical, a 4H-[1,2,4]-oxadiazol-5-one-3-yl radical, nor a 4'-[2,2';6',2'']terpyridinyl radical;
- b) and L¹ is single bond, then L² cannot comprise -N(H)C(=O)C(=O)N(H)- nor -N(H)C(=Q)C(H)CNC(=O)- (where Q is S or O);
- c) and L¹ is other than single bond, then A cannot be quinolin-2-yl-L¹, quinolin-3-yl-L¹, or quinolin-4-yl-L¹;

~~2) when ring A is a fused aryl system, then L¹ must be a single bond;~~

~~3) when ring B is phenyl, ring C is a C₆₋₁₆carbocyclic, L¹ is a single bond, and the compound comprises ring B-OCH₂C(=O)N(H) then ring A cannot be a 2,5-dimethyl-1H-pyrrole-1-yl radical;~~

~~4) ring A cannot be a pyrimidin-2-yl radical when L¹ is -N(H) and ring B is phenyl;~~

~~5) when the compound comprises the formula,~~



where V is =C(H) or =N, and there is a nitrogen of L² bound directly to ring B, then A can not comprise a [1,2,4]-oxadiazol-3-yl radical; and

- 6) the compound is not one of: N-naphthalen-1-yl-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(phenyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-diethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(ethyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(dimethylamino)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl]acetamide, methyl 4-(((3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl)amino]benzoate, ethyl 4-(((3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl)amino]benzoate, 3-(((3-(1H-tetrazol-1-yl)phenyl]

oxy}acetyl)amino]benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-chloro-5-(trifluoromethyl) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acet-amide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.

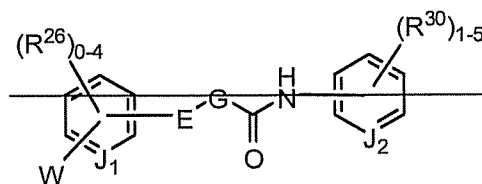
2-12. (cancelled)

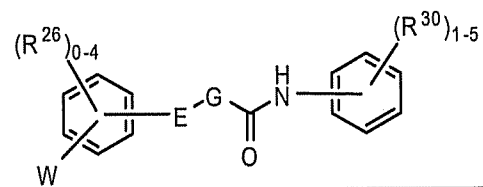
13. (original) The compound according to claim 10, 1, wherein there exists at least one of R³ that is trifluoromethyl.

14. (original) The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L².

15. (currently amended) The compound according to claim 10, wherein each of R³ is independently selected from halogen, trihalomethyl, -OR⁴, -C(=O)R⁴, and optionally substituted C₁₋₆alkyl.

16. (currently amended) A compound for modulating c-Kit activity according to the following Formula: **H**,





II

or a pharmaceutically acceptable salt, thereof, wherein,

W is selected from the following:

each of R^{27} independently selected from halogen, trihalomethyl, $-\text{CN}$, $-\text{NO}_2$, $-\text{OR}^{55}$, $-\text{S(O)}_{0-2}\text{R}^{55}$, $-\text{SO}_2\text{N(R}^{55})\text{R}^{55}$, $-\text{C(=O)N(R}^{55})\text{R}^{55}$, $-\text{C(=NR}^{50})\text{N(R}^{55})\text{R}^{55}$, $-\text{C(=NR}^{50})\text{R}^{55}$, $-\text{N(R}^{55})\text{SO}_2\text{R}^{55}$, $-\text{N(R}^{55})\text{C(O)R}^{55}$, $-\text{NCO}_2\text{R}^{55}$, $-\text{C(=O)R}^{55}$, optionally substituted alkoxy, optionally substituted C_1 -

₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

each Y is independently either =C(H)- or =N-;

Z is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-

E and G are each independently selected from -O-, -S(O)₀₋₂-, -C(R³¹)R³²-, and -N(R³³)-

J₁ and J₂ are each independently =C(H)- or =N-;

R²⁶ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R³⁰ is independently selected from halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl, wherein there exists at least one of R³⁰ that is trihalomethyl; or

~~two adjacent of R²⁶ or two adjacent of R³⁰, together with the annular atoms to which they are attached, can form a five to six membered ring containing up to two heteroatoms and optionally substituted with up to three of R³⁵;~~

R³¹ and R³² are each independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R³³ is selected from -H, optionally substituted lower alkyl, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R⁴⁰ is selected from -H, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two of R⁴⁰, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R⁵⁰ is selected from -H, -CN, -NO₂, -OR⁴⁰, -S(O)₀₋₂R⁴⁰, -CO₂R⁴⁰, optionally substituted C₁₋₆alkyl, optionally substituted C₁₋₆alkenyl, and optionally substituted C₁₋₆alkynyl;

R⁵⁵ is selected from -H, optionally substituted C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl; and

two of R⁵⁵, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

17. (cancelled)

18. (currently amended) The compound according to claim ~~17~~ 16, wherein R³⁰ is selected from halogen, trihalomethyl, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl, wherein there exists at least one of R³⁰ that is trifluoromethyl.

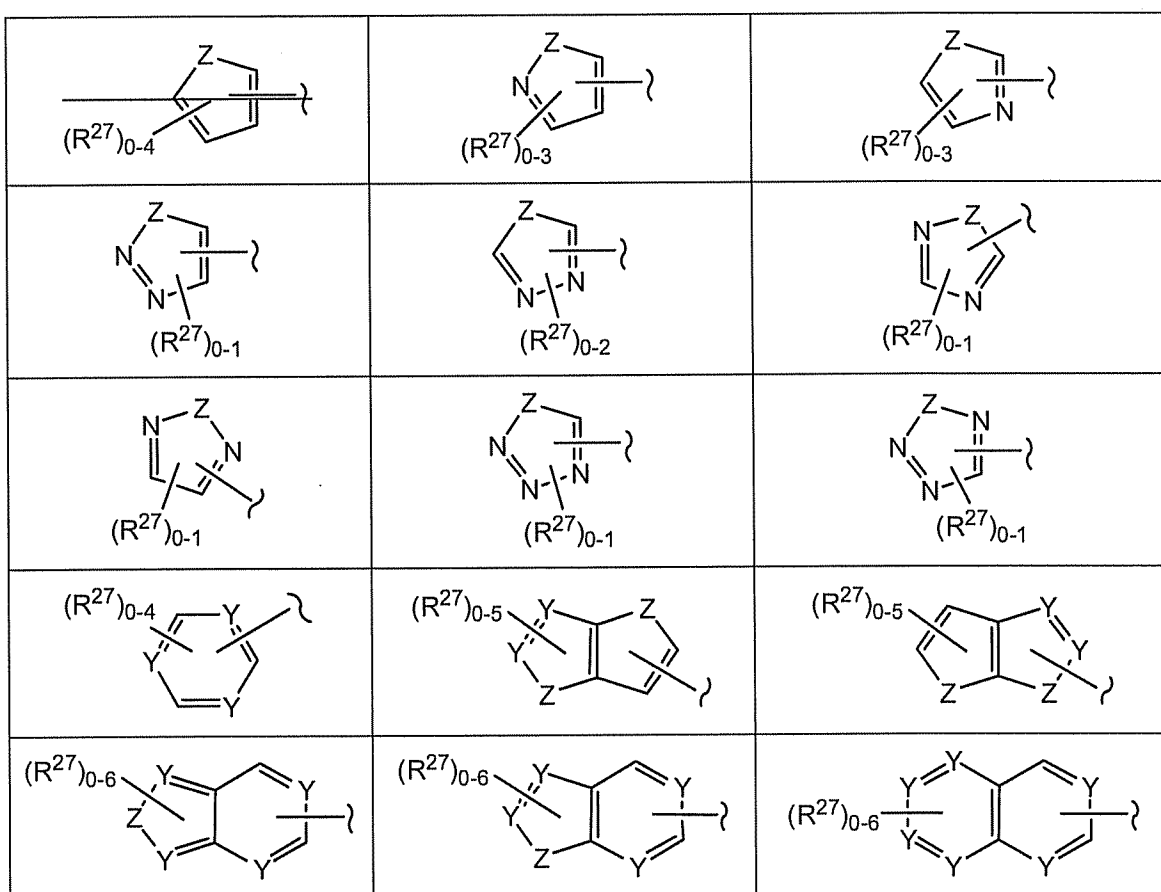
19. (cancelled)

20. (cancelled)

21. (cancelled)

22. (currently amended) The compound according to claim ~~21~~ 16, wherein W is selected from the

following:



and R^{27} is defined as above.

23. (withdrawn from consideration) The compound according to claim 22, wherein E is selected from -O-, -S(O)₀₋₂-, and -NH-; and G is -CH₂-.

24. (withdrawn from consideration) The compound according to claim 22, wherein E is either -CH₂- or -NH-; and G is selected from -O-, -S-, and -NH-.

25. (cancelled)

26. (cancelled)

27. (currently amended) ~~The compound according to claim 1, selected from Table 3:~~

A compound selected from the following Table:

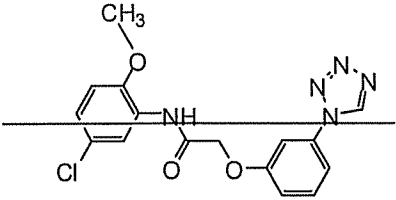
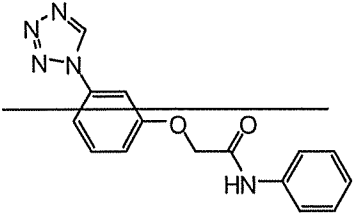
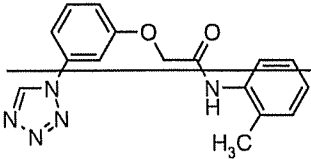
Table 3		
Entry	Name	Structure
1	N-[5-chloro-2-(methyloxy)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
2	N-phenyl 2- {[3-(1H tetrazol 1-yl)phenyl]oxy} acetamide	
3	N-(2-methylphenyl) 2- {[3-(1H tetrazol 1-yl)phenyl]oxy} acetamide	

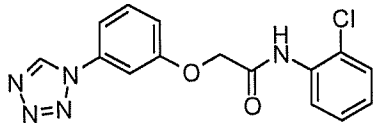
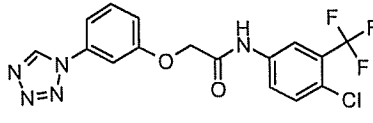
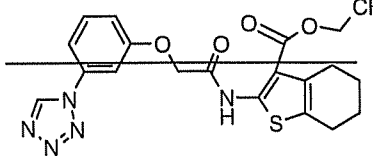
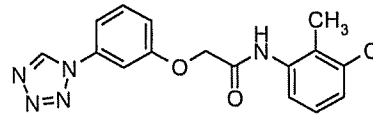
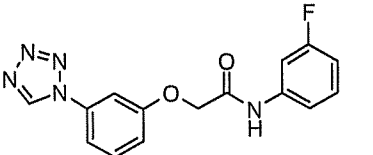
Table 3		
Entry	Name	Structure
4	N-(2-chlorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
5	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
6	ethyl 2-[({[3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl)amino]-4,5,6,7-tetrahydro-1-benzothiophene-3-carboxylate	
7	N-(3-chloro-2-methylphenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
8	N-(3-fluorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	

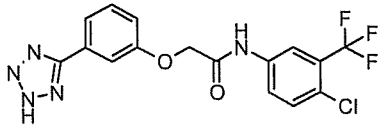
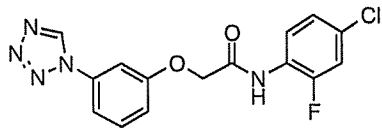
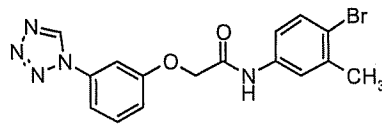
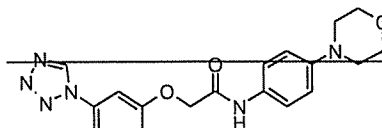
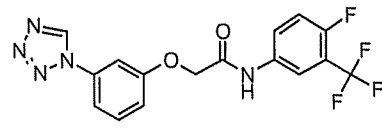
Table 3		
Entry	Name	Structure
9	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2H-tetrazol-5- yl)phenyl]oxy}acetamide	
10	N-(4-chloro-2-fluorophenyl)-2- {[3-(1H- tetrazol-1-yl)phenyl]oxy}acetamide	
11	N-(4-bromo-3-methylphenyl)-2- {[3-(1H- tetrazol-1-yl)phenyl]oxy}acetamide	
12	N-(4-morpholin-4-ylphenyl)-2- {[3-(1H- tetrazol-1-yl)phenyl]oxy}acetamide	
13	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	

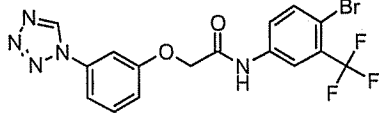
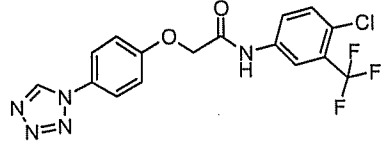
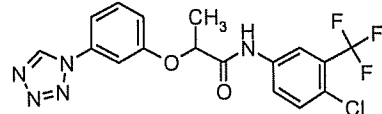
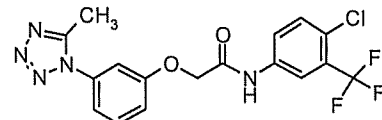
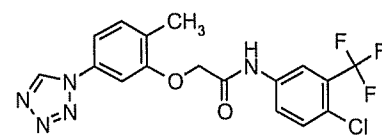
Table 3		
Entry	Name	Structure
14	N-[4-bromo-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy} acetamide	
15	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-tetrazol-1- yl)phenyl]oxy} acetamide	
16	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy} propanamide	
17	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(5-methyl-1H-tetrazol-1- yl)phenyl]oxy} acetamide	
18	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-methyl-5-(1H-tetrazol-1- yl)phenyl]oxy} acetamide	

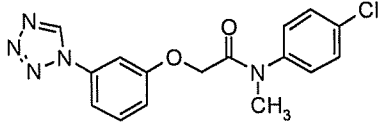
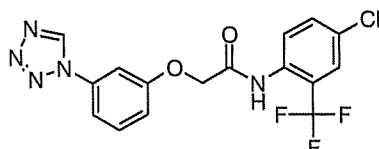
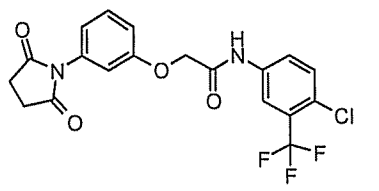
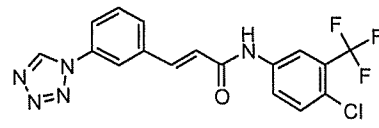
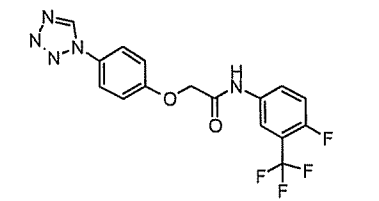
Table 3		
Entry	Name	Structure
19	N-(4-chlorophenyl)-N-methyl-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
20	N-[4-chloro-2-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
21	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(2,5-dioxopyrrolidin-1-yl)phenyl}oxy}acetamide	
22	(2E)-N-[4-chloro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide	
23	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{{4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	

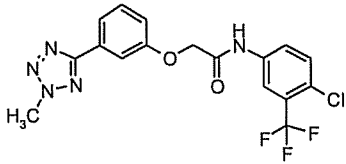
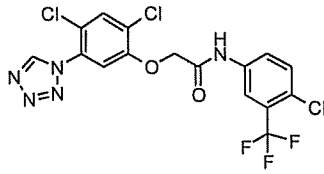
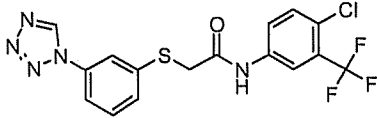
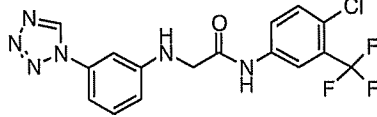
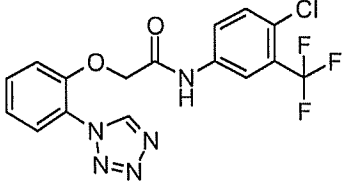
Table 3		
Entry	Name	Structure
24	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2-methyl-2H-tetrazol-5- yl)phenyl]oxy}acetamide	
25	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2,4-dichloro-5-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
26	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]thio}acetamide	
27	N-[4-chloro-3-(trifluoromethyl)phenyl]- N~2~-[3-(1H-tetrazol-1- yl)phenyl]glycinamide	
28	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	

Table 3		
Entry	Name	Structure
29	methyl 1-{3-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-1,2,3-triazole-4-carboxylate	
30	1,1 dimethylethyl {4 [({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]phenyl}carbamate	
31	1,1 dimethylethyl {4 [({[4-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]phenyl}carbamate	
32	N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
33	N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	

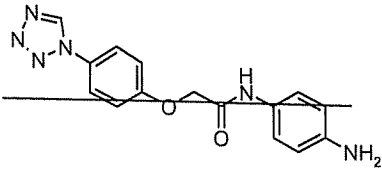
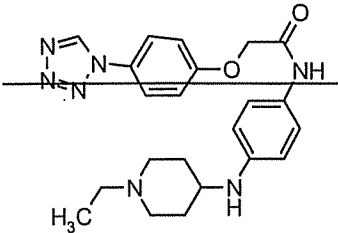
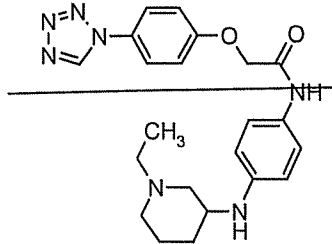
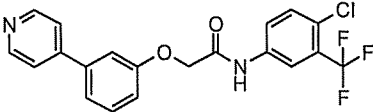
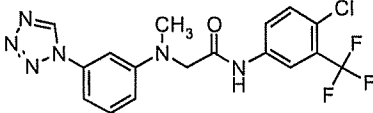
Table 3		
Entry	Name	Structure
34	N-(4-aminophenyl)-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
35	N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
36	N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
37	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-4-ylphenyl)oxy]acetamide	
38	N-[4-chloro-3-(trifluoromethyl)phenyl]-N2-methyl-N2-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	

Table 3		
Entry	Name	Structure
39	N-1,3-benzothiazol-2-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
40	N-quinolin-8-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
41	N-(2,3-dihydro-1,4-benzodioxin-6-yl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
42	N-isoquinolin-5-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
43	N-{3-[(phenylmethyl)oxy]phenyl}-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	

Table 3		
Entry	Name	Structure
44	N-[5-methyl-2-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
45	N-[2,5-bis(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
46	N-(6-fluoro-1,3-benzothiazol-2-yl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
47	methyl 3-[(3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl]amino]benzoate	
48	5-chloro-2-[(3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl]amino]benzamide	

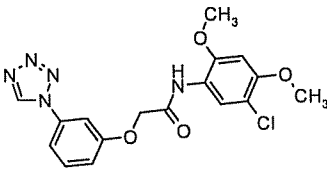
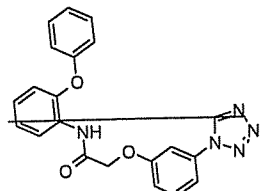
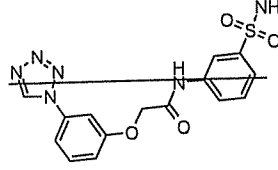
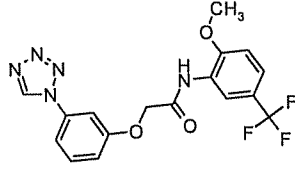
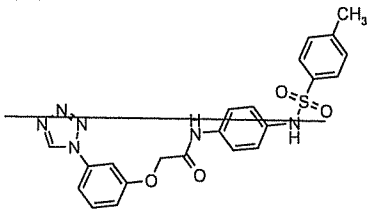
Table 3		
Entry	Name	Structure
49	N-[5-chloro-2,4-bis(methyloxy)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
50	N-[2-(phenyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide	
51	N-[3-(aminosulfonyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide	
52	N-[2-(methyloxy)-5-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide	
53	N-[4-[[[4-methylphenyl)sulfonyl]amino]phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide	

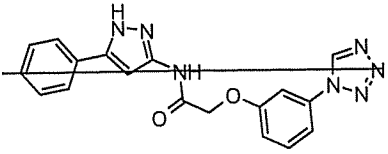
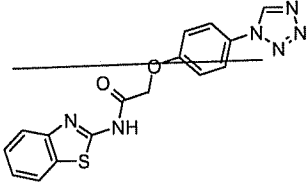
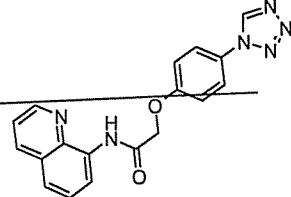
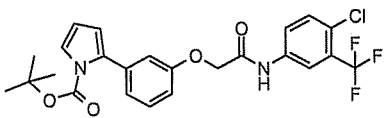
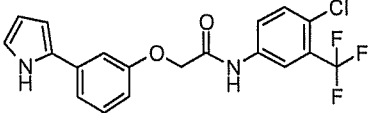
Table 3		
Entry	Name	Structure
54	N-(5-phenyl-1H-pyrazol-3-yl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
55	N-1,3-benzothiazol-2-yl-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
56	N-quinolin-8-yl-2-[[4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
57	1,1-dimethylethyl 2-{3-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-pyrrole-1-carboxylate	
58	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-(1H-pyrrol-2-yl)phenyl]oxy]acetamide	

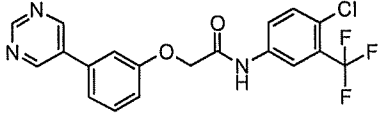
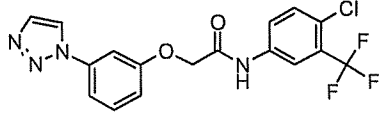
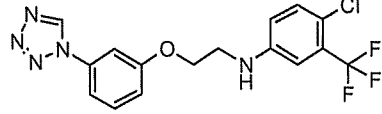
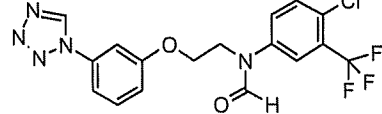
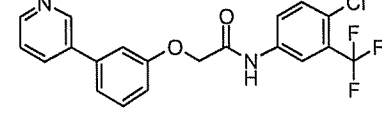
Table 3		
Entry	Name	Structure
59	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyrimidin-5-ylphenyl)oxy]acetamide	
60	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide	
61	4-chloro-N-(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}ethyl)-3-(trifluoromethyl)aniline	
62	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}ethyl)formamide	
63	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-3-ylphenyl)oxy]acetamide	

Table 3		
Entry	Name	Structure
64	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-furan-3-ylphenyl)oxy]acetamide	
65	(2E)-N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide	
66	N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]propanamide	
67	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[6-(1H-tetrazol-1-yl)pyrimidin-4-yl]oxy]acetamide	
68	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-(3,5-dimethylisoxazol-4-yl)phenyl]oxy]acetamide	

Table 3		
Entry	Name	Structure
69	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-quinolin-7-ylphenyl)oxy]acetamide	
70	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-furan-2-ylphenyl)oxy]acetamide	
71	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
72	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-dibenzo[b,d]furan-4-ylphenyl)oxy]acetamide	
73	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	

Table 3		
Entry	Name	Structure
74	N-methyl N-[4-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
75	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[[3-(1H-tetrazol-1-yl)phenyl]methyl]urea	
76	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-methyl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
77	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N~2~-[[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
78	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[[3-(pyridin-2-ylamino)phenyl]oxy]acetamide	

Table 3		
Entry	Name	Structure
79	N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
80	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
81	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
82	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
83	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	

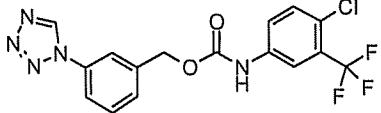
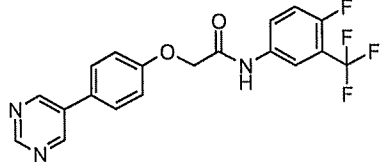
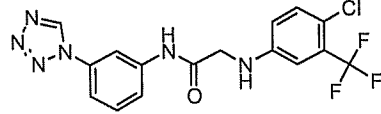
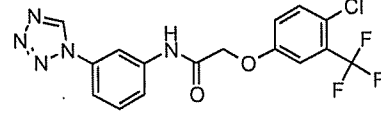
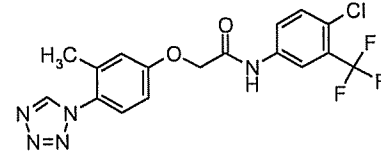
Table 3		
Entry	Name	Structure
84	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
85	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
86	N~2~-[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
87	2-{[4-chloro-3-(trifluoromethyl)phenyl]oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide	
88	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-methyl-4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	

Table 3		
Entry	Name	Structure
89	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-1,2,3-triazol-1- yl)phenyl]oxy}acetamide	
90	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-fluoro-4-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
91	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-fluoro-4-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
92	N-([4-chloro-3- (trifluoromethyl)phenyl]amino)carbonyl)- 3-(1H-tetrazol-1-yl)benzenesulfonamide	
93	N-([4-chloro-3- (trifluoromethyl)phenyl]amino)carbonyl)- N-methyl-3-(1H-tetrazol-1- yl)benzenesulfonamide	

Table 3		
Entry	Name	Structure
94	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
95	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide	
96	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide	
97	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide	
98	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(methyloxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide	

Table 3		
Entry	Name	Structure
99	N-[4-chloro-3-(trifluoromethyl)phenyl]- N~2~-[4-(methoxy)-3-(1H-tetrazol-1- yl)phenyl]glycinamide	
100	N-[4-chloro-3-(trifluoromethyl)phenyl]- N~2~-[4-(1H-tetrazol-1- yl)phenyl]glycinamide	
101	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- (2,3,5,6-tetrafluoro-4-pyrimidin-5- ylphenyl)hydrazinecarboxamide	
102	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(1H-tetrazol-1-yl)phenyl]methyl} urea	
103	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- (4-pyrimidin-5- ylphenyl)hydrazinecarboxamide	

Table 3		
Entry	Name	Structure
104	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
105	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
106	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
107	N-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
108	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	

Table 3		
Entry	Name	Structure
109	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
110	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
111	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
112	1,1-dimethylethyl 2-{4-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-indole-1-carboxylate	
113	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-4-(1H-tetrazol-1-yl)benzenesulfonamide	

Table 3		
Entry	Name	Structure
114	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(2H-tetrazol-5-yl)phenyl]glycinamide	
115	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
116	(3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
117	(3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
118	(3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

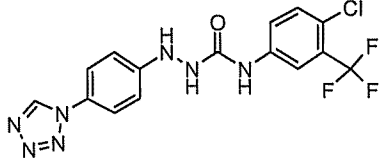
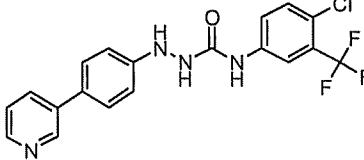
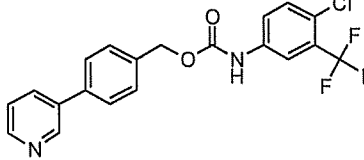
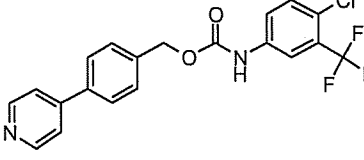
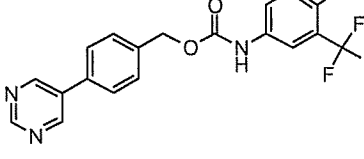
Table 3		
Entry	Name	Structure
119	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
120	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide	
121	(4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
122	(4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
123	(4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3		
Entry	Name	Structure
124	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-4-ylphenyl)methyl]urea	
125	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide	
126	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
127	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
128	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	

Table 3		
Entry	Name	Structure
129	(4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
130	(4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
131	1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
132	1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
133	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	

Table 3		
Entry	Name	Structure
134	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
135	(3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
136	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
137	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
138	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	

Table 3		
Entry	Name	Structure
139	N-{{3-(6-aminopyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
140	N-{{4-(6-aminopyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
141	N-{{3-(2-aminopyrimidin-5-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
142	N-{{4-(2-aminopyrimidin-5-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
143	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyridin-3-ylphenyl)ethyl]urea	

Table 3		
Entry	Name	Structure
144	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyrimidin-5-yl)phenyl]ethylurea	
145	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[4-(1H-indol-2-yl)phenyl]oxy]acetamide	
146	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(isoquinolin-7-yloxy)acetamide	
147	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide	
148	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide	

Table 3		
Entry	Name	Structure
149	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-4-ylphenyl)methyl]urea	
150	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-quinoxalin-6-ylphenyl)methyl]urea	
151	methyl 3-amino-6-(3-{[([4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl]amino}methyl}phenyl)pyrazine-2-carboxylate	
152	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-quinoxalin-6-ylphenyl)methyl]urea	
153	N-{[3-(2-amino-5-methylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3		
Entry	Name	Structure
154	methyl 3-amino-6-(4- {[[[4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl]amino]methyl} phenyl)pyrazine-2-carboxylate	
155	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methoxy)phenyl]carbamate	
156	N-[3-chloro-4-(methoxy)phenyl]-N'--{[3-(1H-tetrazol-1-yl)phenyl]methyl} urea	
157	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(5-hydroxy-1H-tetrazol-1-yl)phenyl]oxy} acetamide	
158	N- {[3-(2-amino-5-chloropyridin-3-yl)phenyl]methyl} -N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3		
Entry	Name	Structure
159	N-{{4-(2-amino-5-chloropyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
160	N-{{3-(6-chloropyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
161	N-{{4-(6-chloropyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
162	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{{4-(pyrimidin-2-yloxy)phenyl}methyl} urea	
163	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-3-(1H-tetrazol-1-yl)benzamide	

Table 3		
Entry	Name	Structure
164	3-amino-6-(3-{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl}amino)methyl}phenyl)-N-[2-(dimethylamino)ethyl]pyrazine-2-carboxamide	
165	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{{3-(6-fluoropyridin-3-yl)phenyl}methyl}urea	
166	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
167	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{{4-(6-fluoropyridin-3-yl)phenyl}methyl}urea	
168	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	

Table 3		
Entry	Name	Structure
169	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(6-methylpyridin-3-yl)phenyl]methyl}urea	
170	N-{[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
171	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-methylpyridin-3-yl)phenyl]methyl}urea	
172	N-{[4-(2-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
173	N-{[3-(2-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3		
Entry	Name	Structure
174	[3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
175	[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
176	[3-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
177	(3-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
178	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(hydroxymethyl)pyridin-3-yl]phenyl} methyl)urea	

Table 3		
Entry	Name	Structure
179	N-{{3-(6-acetylpyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
180	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{{3-(6-cyanopyridin-3-yl)phenyl}methyl}urea	
181	1,1-dimethylethyl (3S)-3-({[3-amino-6-(3-{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl}amino)methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate	
182	3-amino-6-(3-{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl}amino)methyl}phenyl)-N-[(3S)-piperidin-3-yl]pyrazine-2-carboxamide	

Table 3		
Entry	Name	Structure
183	1,1-dimethylethyl (3S)-3-({[3-amino-6-(4-{[([4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl} amino)piperidine-1-carboxylate	
184	3-amino-6-(4-{[([4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl)amino]methyl}phenyl)-N-[(3S)-piperidin-3-yl]pyrazine-2-carboxamide	
185	[3-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
186	N-{[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3		
Entry	Name	Structure
187	[6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
188	[3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
189	[3-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
190	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[5-(methylthio)pyridin-3-yl]phenyl} methyl)urea	
191	[4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3		
Entry	Name	Structure
192	[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
193	[4-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
194	(4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
195	[4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
196	[4-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3		
Entry	Name	Structure
197	[3-(1H-tetrazol-1-yl)phenyl]methyl 1,3-benzothiazol-2-ylcarbamate	
198	[3-(1H-tetrazol-1-yl)phenyl]methyl (5-bromopyridin-2-yl)carbamate	
199	(3-pyridin-3-ylphenyl)methyl (3,5-dimethylphenyl)carbamate	
200	(3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	
201	[4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3		
Entry	Name	Structure
202	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	
203	(4-pyrimidin-5-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
204	(3-pyridin-3-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
205	1,1-dimethylethyl 3-({[3-amino-6-(3-({[4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl)amino]methyl} phenyl)pyrazin-2-yl]carbonyl} amino)piperidine-1-carboxylate	

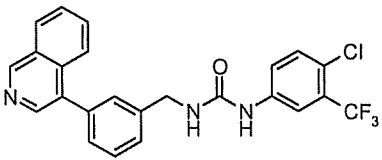
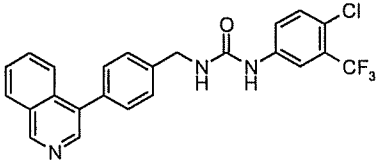
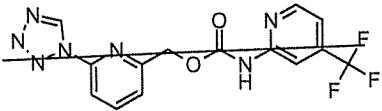
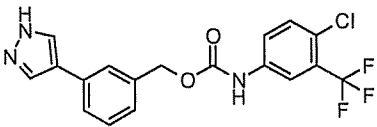
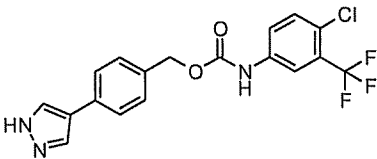
Table 3		
Entry	Name	Structure
206	1,1-dimethylethyl 3-({[3-amino-6-(4- {[(4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl} amino)piperidine-1-carboxylate	
207	3-amino-6-(3-({[(4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl)amino]methyl}phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide	
208	3-amino-6-(4-({[(4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl)amino]methyl}phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide	
209	1,1-dimethylethyl 4-({[3-amino-6-(3- {[(4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl} piperazine-1-carboxylate	

Table 3		
Entry	Name	Structure
210	1,1-dimethylethyl 4-{{3-amino-6-(4- {{{[4-chloro-3-(trifluoromethyl)phenyl]amino} carbonyl]amino}methyl}phenyl)pyrazin-2-yl}carbonyl}piperazine-1-carboxylate	
211	N-({3-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
212	N-({4-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
213	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-pyrazol-4-yl)phenyl]methyl}urea	

Table 3		
Entry	Name	Structure
214	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl} urea	
215	[3-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
216	[4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
217	N-{[3-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
218	N-{[4-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3		
Entry	Name	Structure
219	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
220	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
221	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate	
222	[3-(1H-tetrazol-1-yl)phenyl]methyl [6-(trifluoromethyl)pyridin-2-yl]carbamate	
223	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	

Table 3		
Entry	Name	Structure
224	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-2-yl]phenyl}methyl)urea	
225	[3-(2,6-dimethylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
226	{3-[5-(methyloxy)pyridin-3-yl]phenyl}methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
227	2,3'-bipyridin-6-ylmethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
228	(6-pyrimidin-5-ylpyridin-2-yl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3		
Entry	Name	Structure
229	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-isoquinolin-4-yl)phenyl)methyl]urea	
230	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-isoquinolin-4-yl)phenyl)methyl]urea	
231	[6-(1H-tetrazol-1-yl)pyridin-2-yl)methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	
232	[3-(1H-pyrazol-4-yl)phenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
233	[4-(1H-pyrazol-4-yl)phenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

28. (previously presented) A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.

29. (cancelled)

30. (withdrawn from consideration) A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to claim 1 or a compound selected from N-naphthalen-1-yl-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-(phenyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(3,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,3-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,5-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(3,5-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,6-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,6-diethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2-(ethyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[3-(ethyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,3-dichlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-bromophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-N-[3-(trifluoromethyl)phenyl] acetamide, methyl 4-(((3-(1H-tetrazol-1-yl)phenyl)oxy)acetyl)amino] benzoate, ethyl 4-(((3-(1H-tetrazol-1-yl)phenyl)oxy)acetyl)amino] benzoate, 3-(((3-(1H-tetrazol-1-yl)phenyl)oxy)acetyl)amino] benzoic acid, N-[3-

(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl] oxy} acetamide, N-(4-aminophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide.

31. (withdrawn from consideration) The method according to claim 30, wherein the kinase is c-Kit.
32. (withdrawn from consideration) The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.
33. (withdrawn from consideration) A method of treating diseases or disorders associated with uncontrolled, abnormal, and/or unwanted cellular activities, the method comprising administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in claim 1 or a compound, or a pharmaceutical composition comprising said compound, selected from N-naphthalen-1-yl-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl] -2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethyl-phenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl) acetamide, N-(2-ethyl-phenyl)-2-{{[3-(1H-tetrazol-1-yl) phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyl-oxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}

acetamide, N-[4-(dimethylamino)-phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dichlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-bromophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[[{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino] benzoate, ethyl 4-[[{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino] benzoate, 3-[[{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl] amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl] oxy} acetamide, N-(4-aminophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide.

34. (withdrawn from consideration) A method of screening for modulators of c-Kit, the method comprising combining the compound according to claim 1 or a compound selected from N-naphthalen-1-yl-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl] -2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl) phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{{[3-(1H-tetrazol-1-

yl)phenyl]oxy} acetamide, N-[3-(ethyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dichlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-bromophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[[{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[[{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[[{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl] oxy} acetamide, N-(4-aminophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.

35. (withdrawn from consideration) A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according to claim 1 or a compound selected from N-naphthalen-1-yl-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl] -2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-(2,4,6-

trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl) phenyl}oxy} acetamide, N-(4-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,6-diethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide, N-(2,3-dichlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide, N-(4-bromophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide, N-(2-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[[{3-(1H-tetrazol-1-yl)phenyl}oxy}acetyl]amino] benzoate, ethyl 4-[[{3-(1H-tetrazol-1-yl)phenyl}oxy}acetyl]amino] benzoate, 3-[[{3-(1H-tetrazol-1-yl)phenyl}oxy}acetyl]amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(4H-1,2,4-triazol-4-yl)phenyl}oxy} acetamide, N-(4-chlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl} oxy} acetamide, N-(4-aminophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, and N-(4-acetylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, to a cell or a plurality of cells.